



Titanium: A High Performance Dialect of Java

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<http://www.cs.berkeley.edu/projects/titanium>

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A Little History

- **Most parallel programs are written using explicit parallelism, either:**
 - Message passing with a SPMD model
 - Usually for scientific applications with C++/Fortran
 - Scales easily
 - Shared memory with a thread C or Java
 - Usually for non-scientific applications
 - Easier to program
- **Take the best features of both for Titanium**
 - Builds on ideas in Split-C, AC, and UPC
 - Safer language and more sophisticated implementation

Titanium

- **Take the best features of threads and MPI**
 - global address space like threads (programming)
 - SPMD parallelism like MPI (performance)
 - local/global distinction, i.e., layout matters (performance)
- **Based on Java, a cleaner C++**
 - classes, automatic memory management
 - compiled to C and then assembly (no JVM)
- **Optimizing compiler**
 - communication and memory optimizations
 - synchronization analysis
 - cache and other uniprocessor optimizations

Summary of Features Added to Java

- **Scalable parallelism:**
 - SPMD model of execution with global address space
- **Multidimensional arrays with iterators**
- **Checked Synchronization**
- **Immutable classes**
 - user-definable non-reference types for performance
- **Operator overloading**
- **Zone-based memory management**
- **Libraries**
 - Global communication
 - Distributed arrays
 - Fast bulk I/O

Lecture Outline

- **Language and compiler support for uniprocessor performance**
 - Immutable classes
 - Multidimensional Arrays
 - foreach
- **Language support for parallel computation**
- **Applications and application-level libraries**
- **Summary and future directions**

Java: A Cleaner C++

- **Java is an object-oriented language**
 - classes (no standalone functions) with methods
 - inheritance between classes

- **Documentation on web at java.sun.com**

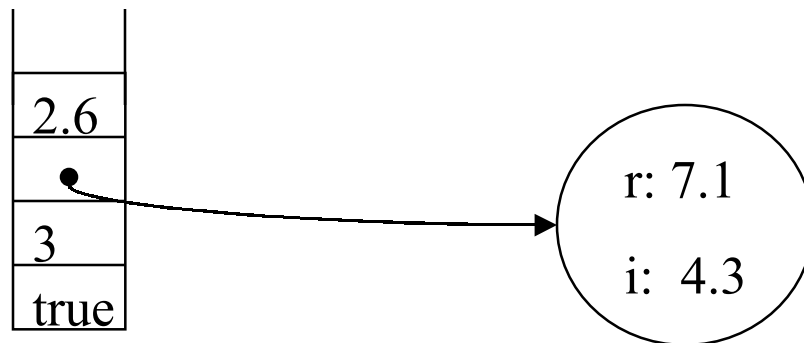
- **Syntax similar to C++**

```
class Hello {  
    public static void main (String [] argv) {  
        System.out.println("Hello, world!");  
    }  
}
```

- **Safe: strongly typed, auto memory management**
- **Titanium is (almost) strict superset**

Java Objects

- **Primitive scalar types: boolean, double, int, etc.**
 - implementations will store these on the program stack
 - access is fast -- comparable to other languages
- **Objects: user-defined and standard library**
 - passed by pointer value (object sharing) into functions
 - has level of indirection (pointer to) implicit
 - simple model, but inefficient for small objects



Java Object Example

```
class Complex {  
    private double real;  
    private double imag;  
    public Complex(double r, double i) {  
        real = r; imag = i; }  
    public Complex operator+(Complex c) {  
        return new Complex(c.real + real,  
                             c.imag + imag); }  
    public double getReal {return real; }  
    public double getImag {return imag; }  
}
```

```
Complex c = new Complex(7.1, 4.3);  
c = c + c;
```

Immutable Classes in Titanium

- **For small objects, would sometimes prefer**
 - to avoid level of indirection
 - pass by value (copying of entire object)
 - especially when immutable -- fields never modified
 - extends the idea of primitive values to user-defined values
- **Titanium introduces immutable classes**
 - all fields are **final** (implicitly)
 - **cannot inherit** from or be inherited by other classes
 - needs to have 0-argument constructor

Example of Immutable Classes

- The immutable complex class nearly the same

new keyword → `immutable` class Complex { *Zero-argument constructor required*
 Complex () {real=0; imag=0; }
 ...
}

Rest unchanged. No assignment to fields outside of constructors.

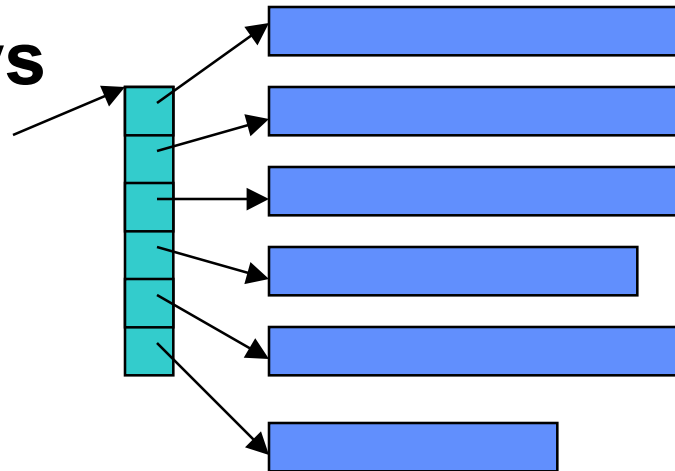
- Use of immutable complex values

```
Complex c1 = new Complex(7.1, 4.3);  
Complex c2 = new Complex(2.5, 9.0);  
c1 = c1 + c2;
```

Similar to structs in C in terms of performance

Arrays in Java

- **Arrays in Java are objects**
- **Only 1D arrays are directly supported**
- **Array bounds are checked**
 - Safe but potentially slow
- **Multidimensional arrays as arrays-of-arrays**
 - General, but slow



Multidimensional Arrays in Titanium

- **New kind of multidimensional array added**
 - Subarrays are supported (unlike Java arrays)
 - Indexed by Points (tuple of ints)
 - Constructed over a set of Points, called Domains
 - RectDomains (rectangular domains) are a special case
 - Points, Domains, RectDomains are immutable classes
- **Support for adaptive meshes and other mesh/grid operations**
 - e.g., can refer to the boundary region of an array

Point, RectDomain, Arrays in General

- **Points specified by a tuple of ints**

```
Point<2> lb = [1, 1];  
Point<2> ub = [10, 20];
```

- **RectDomains given by 3 points:**

- lower bound, upper bound (and stride)

```
RectDomain<2> r = [lb : ub];
```

- **Array declared by # dimensions and type**

```
double [2d] a;
```

- **Array created by passing RectDomain**

```
a = new double [r];
```

Simple Array Example

- Matrix sum in Titanium

```
Point<2> lb = [1,1];
```

```
Point<2> ub = [10,20];
```

```
RectDomain<2> r = [lb,ub];
```

No array allocation here

```
double [2d] a = new double [r];
```

```
double [2d] b = new double [1:10,1:20];
```

```
double [2d] c = new double [lb:ub:[1,1]];
```

Syntactic sugar

```
for (int i = 1; i <= 10; i++)
```

```
    for (int j = 1; j <= 20; j++)
```

```
        c[i,j] = a[i,j] + b[i,j];
```

Optional stride

Naïve MatMul with Titanium Arrays

```
public static void matMul(double [2d] a,  
    double [2d] b, double [2d] c) {  
    int n = c.domain().max()[1]; // square  
    for (int i = 0; i < n; i++) {  
        for (int j = 0; j < n; j++) {  
            for (int k = 0; k < n; k++) {  
                c[i,j] += a[i,k] * b[k,j];  
            }  
        }  
    }  
}
```

Array Performance Issues

- **Array representation is fast, but access methods can be slow, e.g., bounds checking, strides**
- **Compiler optimizes these**
 - common subexpression elimination
 - eliminate (or hoist) bounds checking
 - strength reduce: e.g., naïve code has 1 divide per dimension for each array access
- **Currently +/- 20% of C/Fortran for large loops**
- **Future: small loop and cache optimizations**

Unordered iteration

- All of these optimizations require loop analysis
- Compilers can do this for simple operations, e.g., matrix multiply, but hard in general
- Titanium adds unordered iteration on rectangular domains -- gives user more control

```
foreach (p within r) { ... }
```

- p is a Point new point within the foreach body
- r is a previously-declared RectDomain

Laplacian Example

- Simple example of using arrays and foreach

```
Domain<2> interior = A.domain().shrink(1);
Point<2> dx = [1,0];
Point<2> dy = [0,1];
foreach (p in interior) {
    L[p] = 4*a[p] - a[p+dx] - a[p-dx]
           - a[p+dy] - a[p-dy];
}
```

Better MatMul with Titanium Arrays

```
public static void matMul(double [2d] a,  
    double [2d] b, double [2d] c) {  
    foreach (ij within c.domain()) {  
        double [1d] aRowi = a.slice(1, ij[1]);  
        double [1d] bColj = b.slice(2, ij[2]);  
        foreach (k within aRowi.domain()) {  
            c[ij] += aRowi[k] * bColj[k];  
        }  
    }  
}
```

Current performance: comparable to 3 nested loops in C

Sequential Performance

Ultrasparc:	C/C++/ FORTRAN	Java Arrays	Titanium Arrays	Overhead
DAXPY	1.4s	6.8s	1.5s	7%
3D multigrid	12s		22s	83%
2D multigrid	5.4s		6.2s	15%
MatMul	1.8s		2.2s	22%

Pentium II:	C/C++/ FORTRAN	Java Arrays	Titanium Arrays	Overhead
DAXPY	1.8s		2.3s	27%
3D multigrid	23.0s		20.0s	-13%
2D multigrid	7.3s		5.5s	-25%

Compares to naïve C code; neither compiler does cache blocking (yet).

Lecture Outline

- **Language and compiler support for uniprocessor performance**
- **Language support for parallel computation**
 - **SPMD execution**
 - **Barriers and single**
 - **Explicit Communication**
 - **Implicit Communication (global and local references)**
 - **More on Single**
 - **Synchronized methods and blocks (as in Java)**
- **Applications and application-level libraries**
- **Summary and future directions**

SPMD Execution Model

- **Java programs can be run as Titanium, but the result will be that all processors do all the work**
- **E.g., parallel hello world**

```
class HelloWorld {  
    public static void main (String [] argv) {  
        System.out.println(`Hello from proc ` +  
                             Ti.thisProc());  
    }  
}
```

- **Any non-trivial program will have communication and synchronization**

SPMD Execution Model

- **A common style is compute/communicate**
- **E.g., in each timestep within particle simulation with gravitation attraction**

```
read all particles and compute forces on mine  
Ti.barrier();  
write to my particles using new forces  
Ti.barrier();
```

SPMD Model

- **All processor start together and execute same code, but not in lock-step**
- **Basic control done using**
 - `Ti.numProcs()` total number of processors
 - `Ti.thisProc()` number of executing processor
- **Sometimes they take different branches**

```
if (Ti.thisProc() == 0) { .... do setup .... }  
System.out.println('Hello from ' + Ti.thisProc());  
double [1d] a = new double [Ti.numProcs()];
```

Barriers and Single

- **Common source of bugs is barriers or other global operations inside branches or loops**

`barrier, broadcast, reduction, exchange`

- **A “single” method is one called by all procs**

`public single static void allStep(...)`

- **A “single” variable has same value on all procs**

`int single timestep = 0;`

- **Single annotation on methods (also called “sglobal”) is optional, but useful to understanding compiler messages.**

Explicit Communication: Broadcast

- **Broadcast is a one-to-all communication**

```
broadcast <value> from <processor>
```

- **For example:**

```
int count = 0;  
int allCount = 0;  
if (Ti.thisProc() == 0) count = computeCount();  
allCount = broadcast count from 0;
```

- **The processor number in the broadcast must be single; all constants are single.**
- **The allCount variable could be declared single.**

Example of Data Input

- Same example, but reading from keyboard
- Shows use of Java exceptions

```
int single count = 0;
int allCount = 0;
if (Ti.thisProc() == 0)
    try {
        DataInputStream kb = new DataInputStream(System.in);
        myCount = Integer.valueOf(kb.readLine()).intValue();
    } catch (Exception e) {
        System.err.println("`Illegal Input'");
    }
allCount = myCount from 0;
```

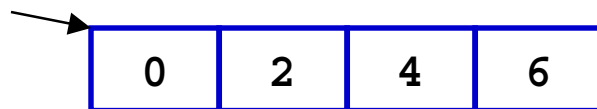
Explicit Communication: Exchange

- **To create shared data structures**
 - each processor builds its own piece
 - pieces are exchanged (for object, just exchange pointers)

- **Exchange primitive in Titanium**

```
int [1d] single allData;  
allData = new int [0:Ti.numProcs()-1];  
allData.exchange(Ti.thisProc()*2);
```

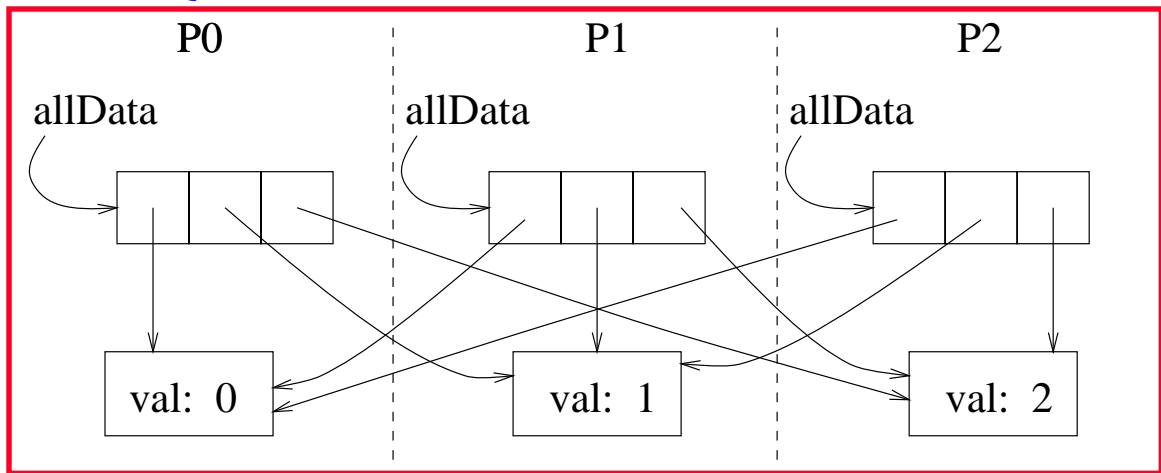
- **E.g., on 4 procs, each will have copy of allData:**



Exchange on Objects

- **More interesting example:**

- ```
class Boxed {
 public Boxed (int j) {
 val = j;
 }
 public in val;
}
```



- ```
Object [1d] single allData;  
allData = new Object [0:Ti.numProcs()-1];  
allData.exchange(new Boxed(Ti.thisProc()));
```

Distributed Data Structures

- **Build distributed data structures with arrays:**

```
RectDomain <1> single allProcs = [0:Ti.numProcs-1];  
RectDomain <1> myParticleDomain = [0:myPartCount-1];  
Particle [1d] single [1d] allParticle =  
    new Particle [allProcs] [1d];  
Particle [1d] myParticle =  
    new Particle [myParticleDomain];  
allParticle.exchange(myParticle);
```

- **Now each processor has array of pointers,
one to each processor's chunk of particles**

More on Single

- **Global synchronization needs to be controlled**
 - if (this processor owns some data) {
 - compute on it
 - barrier
 - }
- **Hence the use of “single” variables in Titanium**
- **If a conditional or loop block contains a barrier, all processors must execute it**
 - conditions in such loops, if statements, etc. must contain only single variables

Single Variable Example

- **Barriers and single in N-body Simulation**

```
class ParticleSim {  
    public static void main (String [] argv) {  
        int single allTimestep = 0;  
        int single allEndTime = 100;  
        for (; allTimestep < allEndTime; allTimestep++){  
            read all particles and compute forces on mine  
            Ti.barrier();  
            write to my particles using new forces  
            Ti.barrier();  
        }  
    }  
}
```

- **Single methods inferred; see David Gay's work**

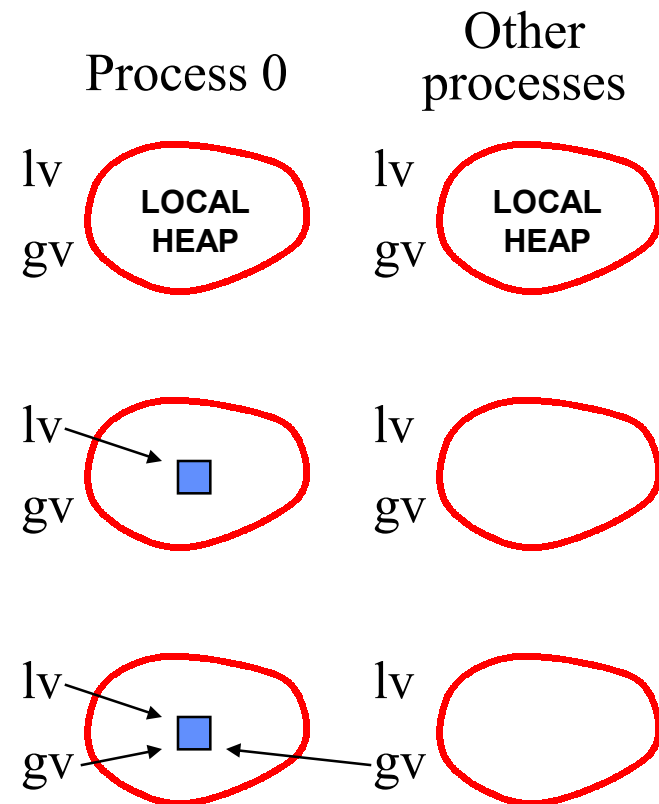
Use of Global / Local

- **As seen, references (pointers) may be remote**
 - easy to port shared-memory programs
- **Global pointers are more expensive than local**
 - True even when data is on the same processor
 - Use **local** declarations in critical sections
- **Costs of global:**
 - space (processor number + memory address)
 - dereference time (check to see if local)
- **May declare references as local**

Global Address Space

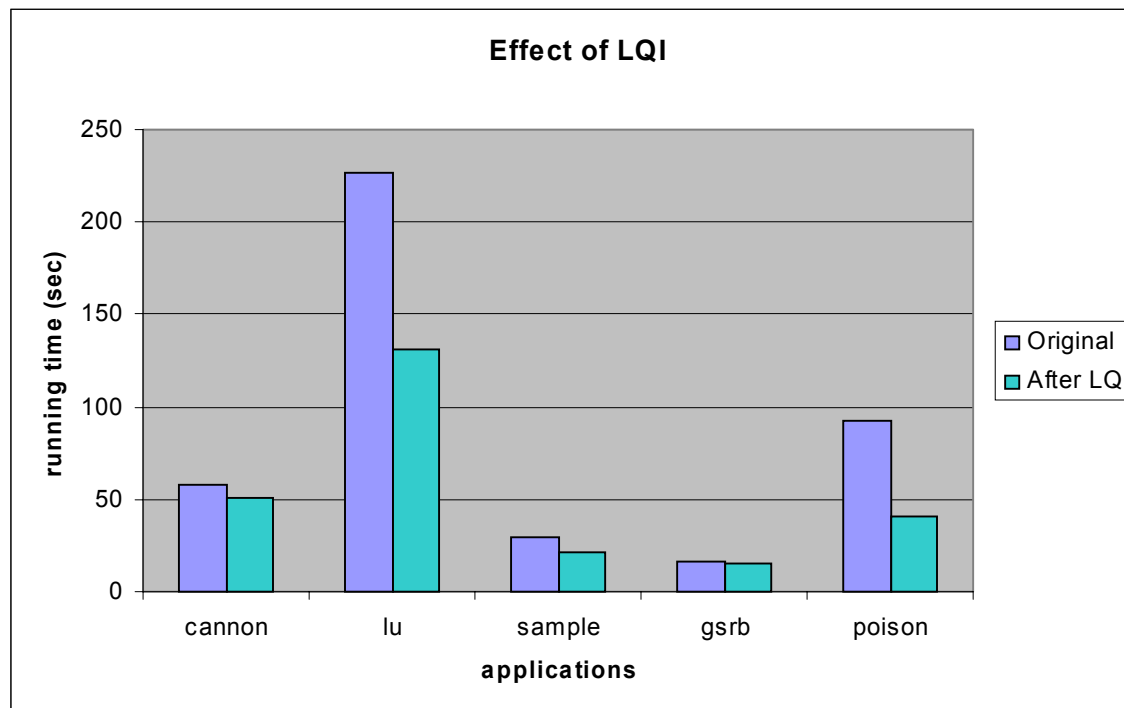
- Processes allocate locally
- References can be passed to other processes

```
Class C { int val;.. }  
C gv;      // global pointer  
C local lv; // local pointer  
  
if (thisProc() == 0) {  
    lv = new C();  
}  
gv = broadcast lv from 0;  
gv.val = ..; // full  
.. = gv.val; // functionality
```



Local Pointer Analysis

- **Compiler can infer many uses of local**
 - See Liblit's work on Local Qualification Inference



- **Data structures must be well partitioned**

Region-Based Memory Management

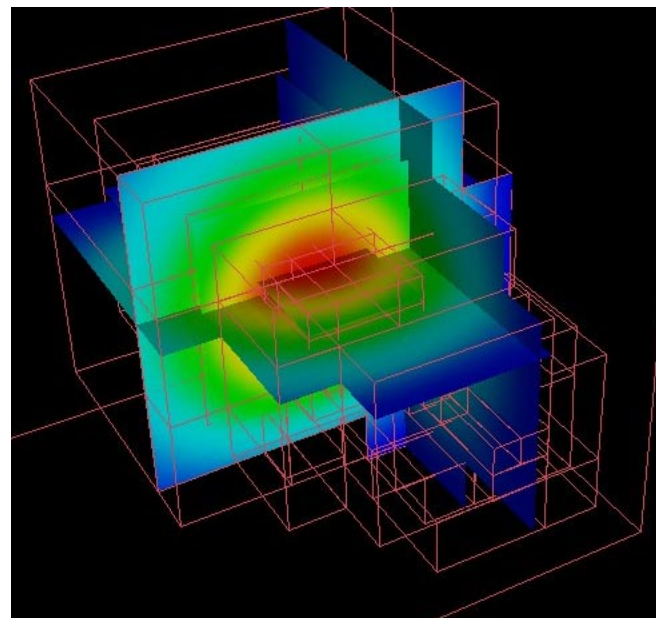
```
PrivateRegion r = new PrivateRegion();
For (int j = 0; j < 10; j++) {
    int[] x = new ( r ) int[j + 1];
    work(j, x);
}
try { r.delete; }
catch (RegionInUse oops) {
    system.out.println("failed to delete");
}
}
```

Lecture Outline

- **Language and compiler support for uniprocessor performance**
- **Language support for parallel computation**
- **Applications and application-level libraries**
 - AMR overview
 - AMR and uniform grid algorithms in Titanium
 - Several smaller benchmarks
 - MatMul, LU, FFT, Join, Sort, EM3d
 - Library interfaces
 - PETSc, Metis,
- **Summary and future directions**

Block-Structured AMR

- **Algorithms for many rectangular, grid-based computations are**
 - communication intensive
 - memory intensive
- **AMR makes these harder**
 - more small messages
 - more complex data structures
 - most of the programming effort is debugging the boundary cases
 - locality and load balance trade-off is hard



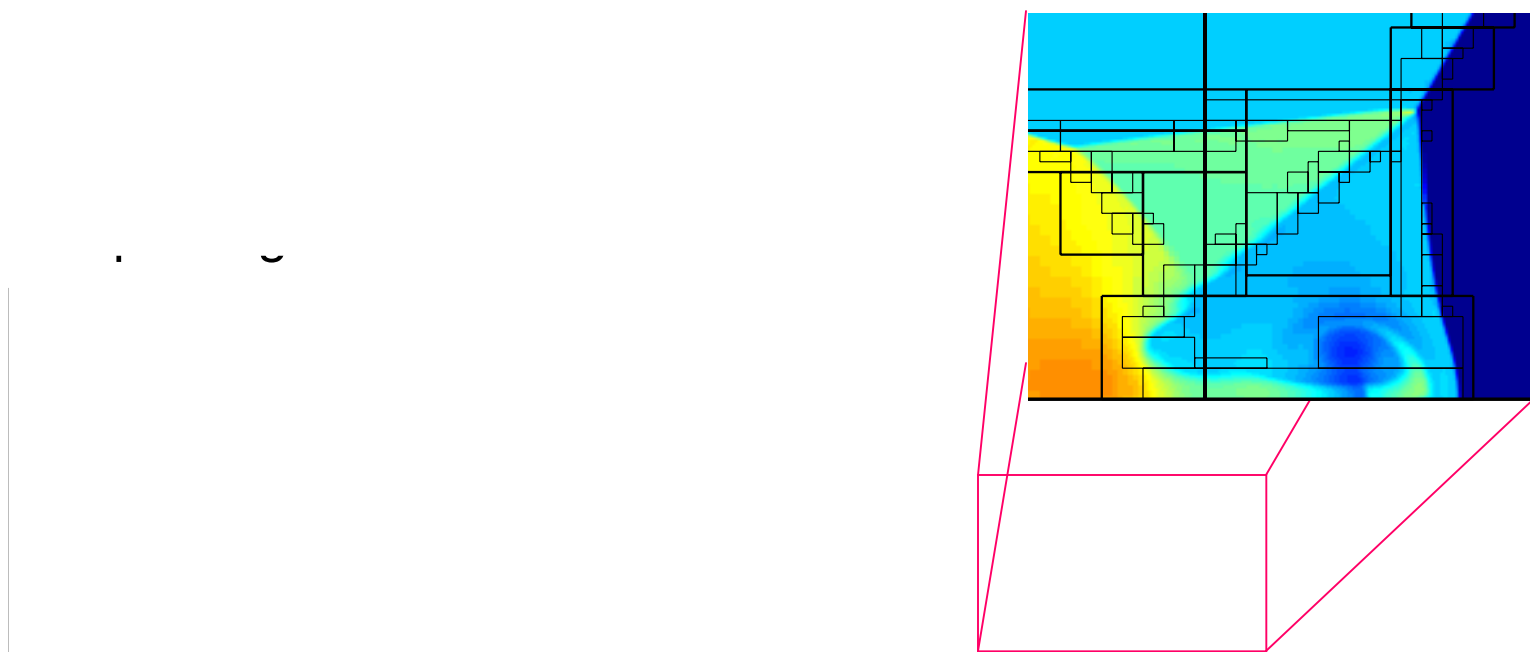
Algorithms for AMR

- **Existing algorithms in Titanium**
 - 3D AMR Poisson solver
 - 3D AMR Gas dynamics
 - Domain-decomposition MLC Poisson
- **Under development**
 - Self-gravitating gas dynamics (3D AMR)
 - For stellar collapse, etc.
 - Immersed boundary method (3D, non-adaptive)
 - Peskin and MacQueen's method for heart model, etc.
 - Embedded boundaries
 - Simulation of bio-MEMs devices and cellular level modeling
 - Project Idea:
 - Multiblock Java code with self-scheduling. Contact me, [yelick@cs.](mailto:yelick@cs.berkeley.edu)
 - Evaluation of and proposal for general domains.
- **All joint with Colella's group at LBNL**

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3D AMR Gas Dynamics



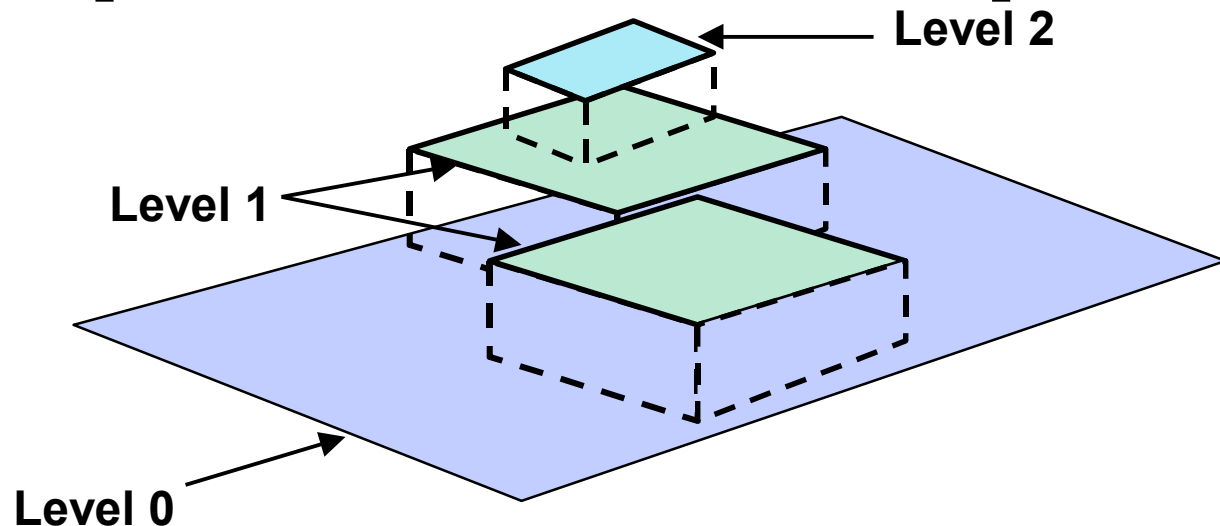
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3D AMR Poisson

- **Poisson Solver [Semenzato, Pike, Colella]**

- finite domain
- variable coefficients
- multigrid across levels



- **Currently synthetic grids, no grid generation**
- **Under construction**
 - reengineered to interface with hyperbolic solver
 - including mesh generation

MLC for Finite-Differences

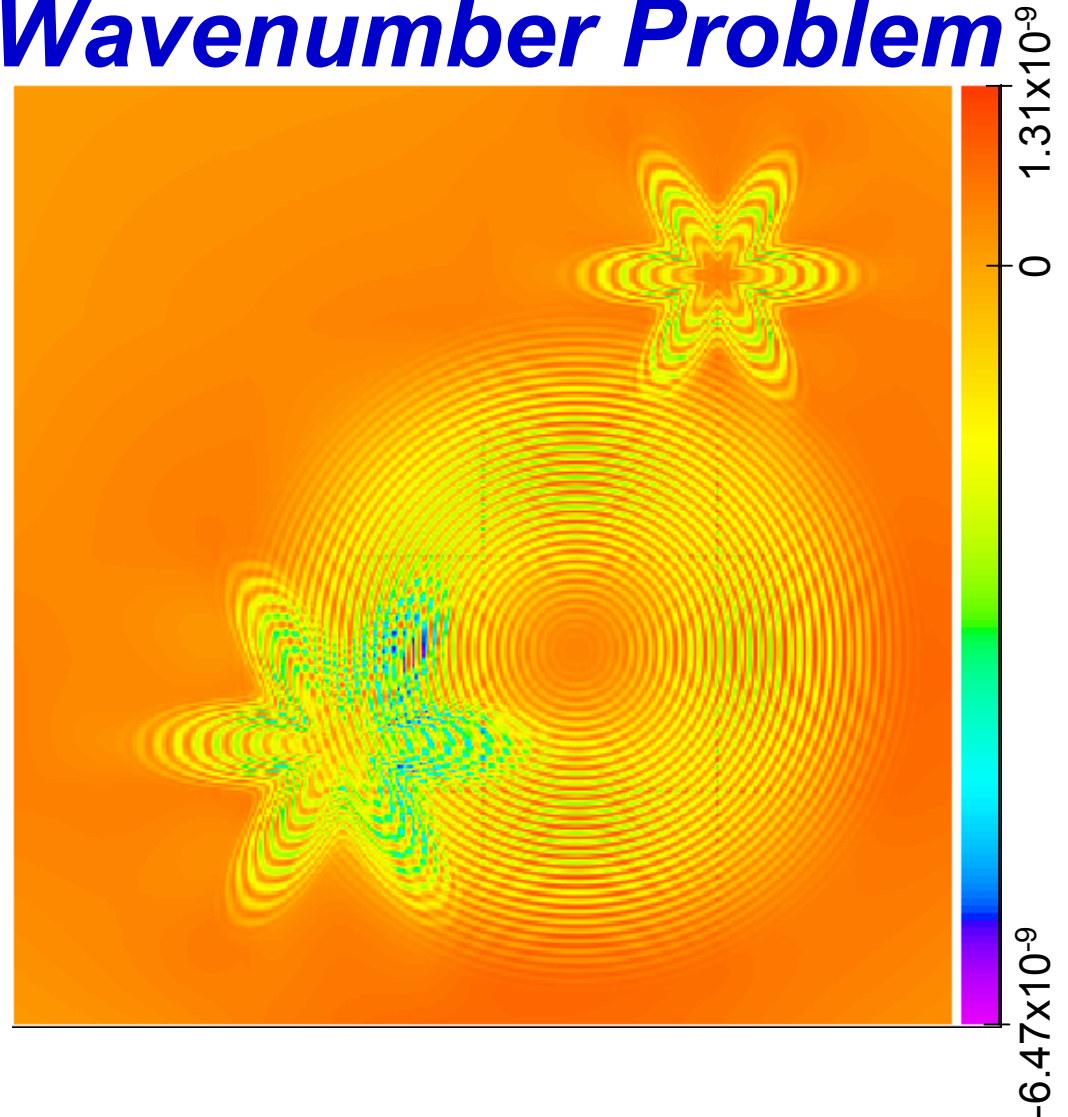
- **Poisson solver with infinite domains [Colella, Balls]**
 - Uses a Method of Local Corrections (MLC)
 - Currently non-adaptive and 2D
 - Supports only constant coefficients
- **Uses 2-level, domain decomposition approach**
 - Fine-grid solutions are computed in parallel
 - Information transferred to a coarse-grid and solved serially
 - Fine-grid solutions is computed using boundary conditions from the coarse grid
- **Future work includes 3D Adaptive version**

MLC for Finite-Differences

- **Features of the method**
 - Solution is still second-order accurate
 - Accuracy depends only weakly on the coarse-grid spacing
- **Scalability**
 - No communication during fine-grid solves
 - Single communication step (global all-to-all)
 - coarse grid work is serial (replicated), but relatively small
- **Future work: extend to 3D and adaptive meshes**
 - Project idea: extension to 3D: see Greg Balls, gballs@cs

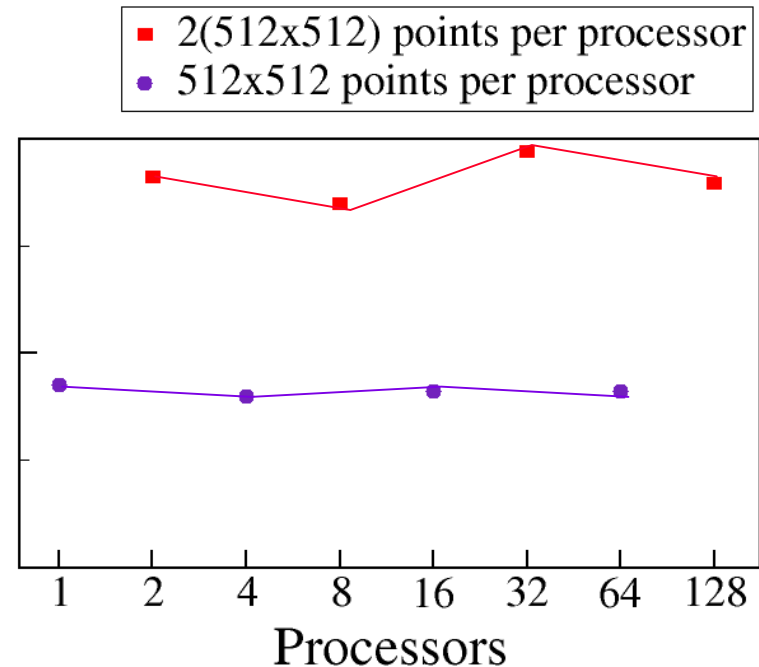
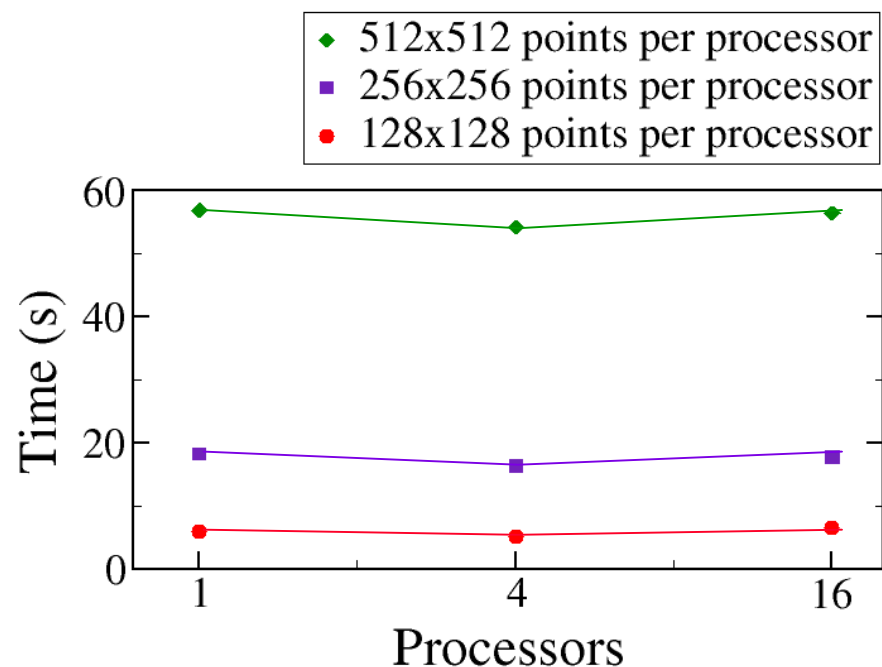
Error on High-Wavenumber Problem

- **Charge is**
 - 1 charge of concentric waves
 - 2 star-shaped charges.
- **Largest error is where the charge is changing rapidly. Note:**
 - discretization error
 - faint decomposition error
- **Run on 16 procs**



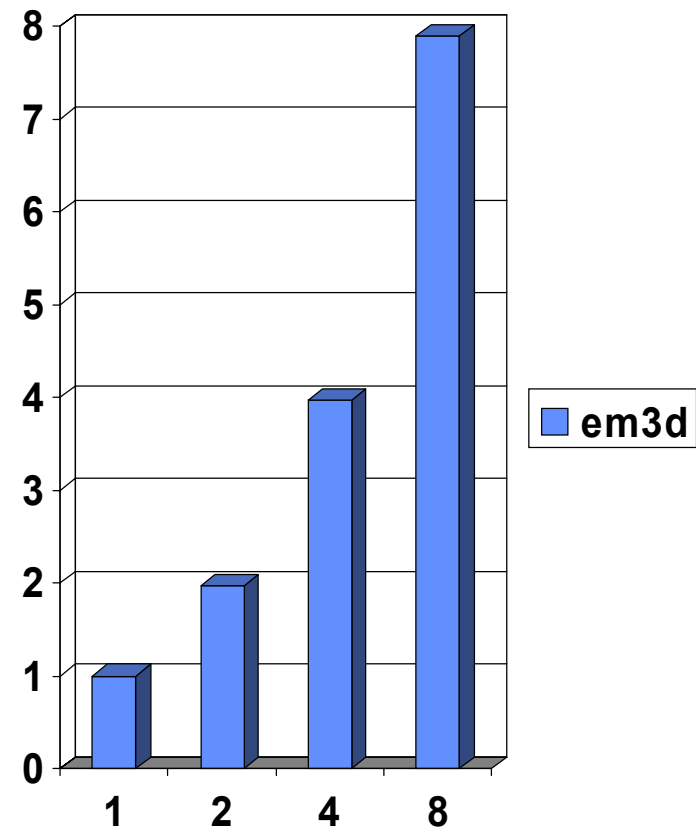
Scalable Poisson Solver (MLC)

- Communication performance is low ($< 5\%$)
- Scaled speedup experiments are nearly ideal (flat)



Unstructured Mesh Kernel

- **EM3D: Relaxation on a 3D unstructured mesh**
- **Speedup on Ultrasparc SMP**
- **Simple kernel: mesh not partitioned.**



Calling Other Languages

- **We have built interfaces to**
 - PETSc : scientific library for finite element applications
 - Metis: graph partitioning library
- **Two issues with cross-language calls**
 - accessing Titanium data structures (arrays) from C
 - possible because Titanium arrays have same format on inside
 - having a common message layer
 - Titanium is built on lightweight communication

Lecture Outline

- **Language and compiler support for uniprocessor performance**
- **Language support for parallel computation**
- **Applications and application-level libraries**
- **Summary and future directions**
 - Implementation

Implementation

- **Strategy**
 - Titanium into C
 - Solaris or Posix threads for SMPs
 - Lightweight communication for MPPs/Clusters
- **Status: Titanium runs on**
 - Solaris or Linux SMPs and uniprocessors
 - Berkeley NOW
 - SDSC Tera, SP2, T3E (NERSC and NPACI)
 - SP3 (and IBM SP Power3) port underway

Titanium Summary

- **Performance**
 - close to C/FORTRAN + MPI on limited class of problems
- **Portability**
 - develop on uniprocessor, then SMP, then MPP/Cluster
- **Safety**
 - as safe as Java, extended to parallel framework
- **Expressiveness**
 - easier than MPI, harder than threads
- **Compatibility, interoperability, etc.**
 - no gratuitous departures from Java standard

Using Titanium

- **On machines in the CS Division**

- `/srs/titanium/*/bin/tcbuild file.ti`

- Solaris 2.6 and Linux supported; need to mount this filesystem

- **On NERSC t3e use:**

- `/u/mp215/miyamoto/tc-1.44/tcbuild/tcbuild file.ti`

- **On SP2 contact:** `cjlin@cs.berkeley.edu`

- **For documentation, source code, see the home page**

- <http://www.cs.berkeley.edu/projects/titanium>

- **Documentation includes**

- Language reference, terse but complete
 - Tutorial, incomplete

- **For problems or questions:**

- `titanium-group@cs.berkeley.edu`

Future Plans

- **Improved compiler optimizations for scalar code**
 - large loops are currently +/- 20% of Fortran
 - working on small loop performance
- **Packaged solvers written in Titanium**
 - Elliptic and hyperbolic solvers, both regular and adaptive
- **New application collaboration**
 - Peskin and McQueen (NYU) with Colella (LBNL)
 - Immersed boundary method, currently use for heart simulation, platelet coagulation, and others



Backup Slides

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Other Language Extensions

Java extensions for expressiveness & performance

- **Operator overloading**
- **Zone-based memory management**
- **Foreign function interface**

The following is not yet implemented in the compiler

- **Parameterized types (aka templates)**

Consistency Model

- **Titanium adopts the Java memory consistency model**
- **Roughly: Access to shared variables that are not synchronized have undefined behavior.**
- **Use synchronization to control access to shared variables.**
 - barriers
 - synchronized methods and blocks